

10/784,673

Connecting via Winsock to STN

Welcome to STN International! Enter x:x

LOGINID:sssptal611txm

PASSWORD:

* * * * * RECONNECTED TO STN INTERNATIONAL * * * * *
SESSION RESUMED IN FILE 'REGISTRY' AT 18:20:53 ON 04 MAY 2005
FILE 'REGISTRY' ENTERED AT 18:20:53 ON 04 MAY 2005
COPYRIGHT (C) 2005 American Chemical Society (ACS)

=> s (nc4-c6 or n2c3-c6 or ncnc2-c6 or n3c2-c6)es
MISSING OPERATOR

=> s (nc4-c6 or n2c3-c6 or ncnc2-c6 or n3c2-c6)/es
833648 NC4-C6/ES
53398 N2C3-C6/ES
242909 NCNC2-C6/ES
41176 N3C2-C6/ES

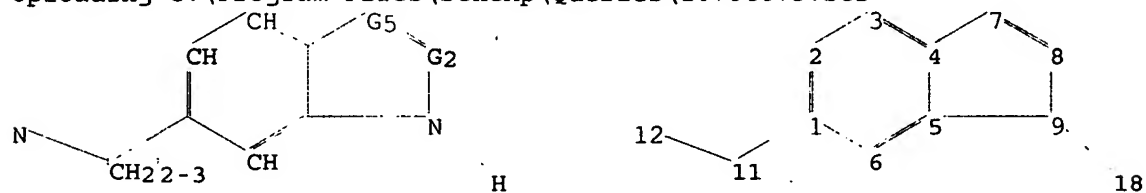
SYSTEM LIMITS EXCEEDED - SEARCH ENDED

The search profile you entered was too complex or gave too many answers. Simplify or subdivide the query and try again. If you have exceeded the answer limit, enter DELETE HISTORY at an arrow prompt (=>) to remove all previous answers sets and begin at L1. Use the SAVE command to store any important profiles or answer sets before using DELETE HISTORY.

=> delete history
DELETE ALL L# ITEMS? (Y)/N:.
ALL L# ITEMS DELETED

=> s (nc4-c6 or n2c3-c6 or ncnc2-c6 or n3c2-c6)/es
833648 NC4-C6/ES
53398 N2C3-C6/ES
242909 NCNC2-C6/ES
41176 N3C2-C6/ES
L1 - 1163389 (NC4-C6 OR N2C3-C6 OR NCNC2-C6 OR N3C2-C6)/ES

=>
Uploading C:\Program Files\Stnexp\Queries\10784673.str



chain nodes :
11 18
ring nodes :
1 2 3 4 5 6 7 8 9
ring/chain nodes :
12
chain bonds :

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1-11 9-18 11-12
ring bonds :
1-2 1-6 2-3 3-4 4-5 4-7 5-6 5-9 7-8 8-9
exact/norm bonds :
1-11 4-7 5-9 7-8 8-9 9-18 11-12
normalized bonds :
1-2 1-6 2-3 3-4 4-5 5-6

G1

G2:N,CH

G5:C,N

Match level :

1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:Atom 8:Atom 9:Atom 11:CLASS
12:CLASS 18:CLASS

L2 STRUCTURE UPLOADED

=> s 12 subset = 11 sample

SAMPLE SUBSET SEARCH INITIATED 18:23:24 FILE 'REGISTRY'

SAMPLE SUBSET SCREEN SEARCH COMPLETED - 43320 TO ITERATE

2.3% PROCESSED 1000 ITERATIONS
INCOMPLETE SEARCH (SYSTEM LIMIT EXCEEDED)
SEARCH TIME: 00.00.01

0 ANSWERS

PROJECTIONS (WITHIN SPECIFIED SUBSET):

ONLINE **INCOMPLETE**

PROJECTED ITERATIONS (WITHIN SPECIFIED SUBSET):

853978 TO 878822

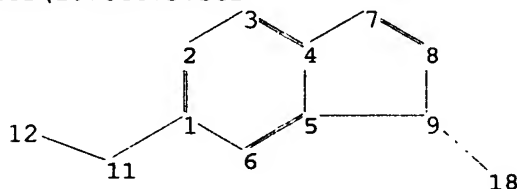
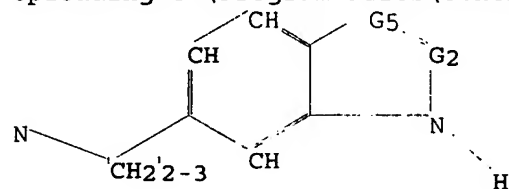
PROJECTED ANSWERS (WITHIN SPECIFIED SUBSET):

0 TO 0

L3 0 SEA SUB=L1 SSS SAM L2

=>

Uploading C:\Program Files\Stnexp\Queries\10784673.str



chain nodes :

11 18

ring nodes :

1 2 3 4 5 6 7 8 9

ring/chain nodes :

12

chain bonds :

1-11 9-18 11-12

ring bonds :

1-2 1-6 2-3 3-4 4-5 4-7 5-6 5-9 7-8 8-9

exact/norm bonds :

1-11 4-7 5-9 7-8 8-9 9-18 11-12

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normalized bonds :

1-2 1-6 2-3 3-4 4-5 5-6

G1

G2:N,CH

G5:C,N

Match level :

1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:Atom 8:Atom 9:Atom 11:CLASS
12:CLASS 18:CLASS

L4 STRUCTURE UPLOADED

=> s l4 subset = l1 sample

SAMPLE SUBSET SEARCH INITIATED 18:24:51 FILE 'REGISTRY'

SAMPLE SUBSET SCREEN SEARCH COMPLETED - 43320 TO ITERATE

2.3% PROCESSED 1000 ITERATIONS
INCOMPLETE SEARCH (SYSTEM LIMIT EXCEEDED)
SEARCH TIME: 00.00.01

0 ANSWERS

PROJECTIONS (WITHIN SPECIFIED SUBSET):

ONLINE **INCOMPLETE**

PROJECTED ITERATIONS (WITHIN SPECIFIED SUBSET):

853978 TO 878822

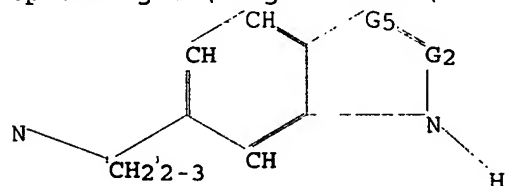
PROJECTED ANSWERS (WITHIN SPECIFIED SUBSET):

0 TO 0

L5 0 SEA SUB=L1 SSS SAM L4

=>

Uploading C:\Program Files\Stnexp\Queries\10784673.str



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G2:N,CH

G5:C,N

Match level :

1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:Atom 8:Atom 9:Atom 11:CLASS
12:CLASS 18:CLASS

L6 STRUCTURE UPLOADED

=> s l6 subset = l1 sample

SAMPLE SUBSET SEARCH INITIATED 18:27:09 FILE 'REGISTRY'

SAMPLE SUBSET SCREEN SEARCH COMPLETED - 14401 TO ITERATE

6.9% PROCESSED 1000 ITERATIONS 0 ANSWERS
INCOMPLETE SEARCH (SYSTEM LIMIT EXCEEDED)
SEARCH TIME: 00.00.01

PROJECTIONS (WITHIN SPECIFIED SUBSET): ONLINE **COMPLETE**
PROJECTED ITERATIONS (WITHIN SPECIFIED SUBSET): 280833 TO 295207
PROJECTED ANSWERS (WITHIN SPECIFIED SUBSET): 0 TO 0

L7 0 SEA SUB=L1 SSS SAM L6

=> s l6 subset = l1 full

FULL SUBSET SEARCH INITIATED 18:27:26 FILE 'REGISTRY'

FULL SUBSET SCREEN SEARCH COMPLETED - 285210 TO ITERATE

100.0% PROCESSED 285210 ITERATIONS 151 ANSWERS
SEARCH TIME: 00.00.02

L8 151 SEA SUB=L1 SSS FUL L6

=> s c12h16n2 and l8

1947 C12H16N2

L9 1 C12H16N2 AND L8

=> d ide cbib pi

L9 ANSWER 1 OF 1 REGISTRY COPYRIGHT 2005 ACS on STN

RN 321744-76-7 REGISTRY

ED Entered STN: 14 Feb 2001

CN 1H-Indole-6-ethanamine, N,N-dimethyl- (9CI) (CA INDEX NAME)

OTHER NAMES:

CN 6-[2-(Dimethylamino)ethyl]-1H-indole

FS 3D CONCORD

MF C12 H16 N2

SR CA

LC STN Files: CA, CAPLUS, USPATFULL

Me₂N-CH₂-CH₂

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

2 REFERENCES IN FILE CA (1907 TO DATE)
2 REFERENCES IN FILE CAPLUS (1907 TO DATE)

REFERENCE 1: 140:303530 Preparation of indole and indazole derivatives for the treatment of migraine. Edwards, Louise; Isaac, Methvin; Maddaford, Shawn; Slassi, Abdelmalik; Xin, Tao (NPS Allelix Biopharmaceuticals, Inc., Can.). U.S. US 6716837 B1 20040406, 29 pp., Cont.-in-part of U.S. Ser. No. 354,091, abandoned. (English). CODEN: USXXAM. APPLICATION: US 2000-709579 20001113. PRIORITY: US 1999-354091 19990715.
PATENT NO. KIND DATE APPLICATION NO. DATE

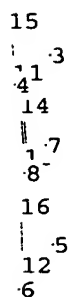
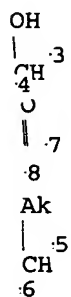
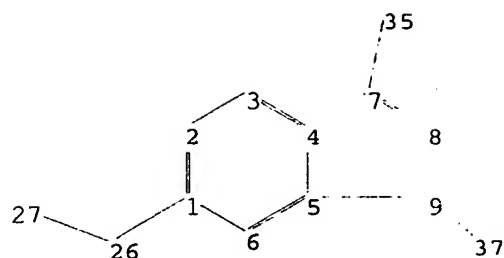
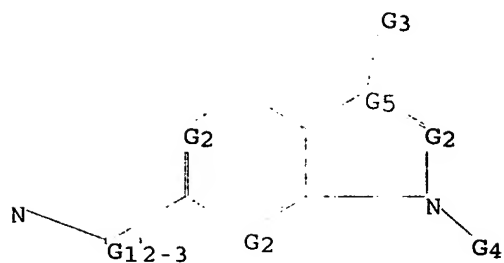
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PI	US 6716837	B1	20040406	US 2000-709579	20001113
	US 2004167158	A1	20040826	US 2004-784673	20040224

REFERENCE 2: 134:131423 Preparation of aminoalkylindoles and analogs as 5-HT_{1D} receptor ligands. Edwards, Louise; Isaac, Methvin; Maddaford, Shawn; Slassi, Abdelmalik; Xin, Tao (NPS Allelix Corp., Can.). PCT Int. Appl. WO 2001005758 A2 20010125, 71 pp. DESIGNATED STATES: W: AE, AG, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BY, BZ, CA, CH, CN, CR, CU, CZ, DE, DK, DM, DZ, EE, ES, FI, GB, GD, GE, GH, GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MA, MD, MG, MK, MN, MW, MX, MZ, NO, NZ, PL, PT, RO, RU, SD, SE, SG, SI, SK, SL, TJ, TM, TR, TT, TZ, UA, UG, US, UZ, VN, YU, ZA, ZW, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM; RW: AT, BE, BF, BJ, CF, CG, CH, CI, CM, CY, DE, DK, ES, FI, FR, GA, GB, GR, IE, IT, LU, MC, ML, MR, NE, NL, PT, SE, SN, TD, TG. (English). CODEN: PIXXD2. APPLICATION: WO 2000-CA831 20000714. PRIORITY: US 1999-354091 19990715.
PATENT NO. KIND DATE APPLICATION NO. DATE

	PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
PI	WO 2001005758	A2	20010125	WO 2000-CA831	20000714
	WO 2001005758	A3	20010719		
	W:				
	AE, AG, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BY, BZ, CA, CH, CN, CR, CU, CZ, DE, DK, DM, DZ, EE, ES, FI, GB, GD, GE, GH, GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MA, MD, MG, MK, MN, MW, MX, MZ, NO, NZ, PL, PT, RO, RU, SD, SE, SG, SI, SK, SL, TJ, TM, TR, TT, TZ, UA, UG, US, UZ, VN, YU, ZA, ZW, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM				
	RW:				
	GH, GM, KE, LS, MW, MZ, SD, SL, SZ, TZ, UG, ZW, AT, BE, CH, CY, DE, DK, ES, FI, FR, GB, GR, IE, IT, LU, MC, NL, PT, SE, BF, BJ, CF, CG, CI, CM, GA, GN, GW, ML, MR, NE, SN, TD, TG				
	CA 2378047	AA	20010125	CA 2000-2378047	20000714
	EP 1196380	A2	20020417	EP 2000-945511	20000714
	R:				
	AT, BE, CH, DE, DK, ES, FR, GB, GR, IT, LI, LU, NL, SE, MC, PT, IE, SI, LT, LV, FI, RO				
	JP 2003505369	T2	20030212	JP 2001-511419	20000714

=>

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Cb⁹32⁹

chain nodes :

10 11 12 13 14 15 16 26 32 35 37

ring nodes :

1 2 3 4 5 6 7 8 9

ring/chain nodes :

27

chain bonds :

1-26 7-35 9-37 11-15 12-16 13-14 26-27

ring bonds :

1-2 1-6 2-3 3-4 4-5 4-7 5-6 5-9 7-8 8-9

exact/norm bonds :

1-2 1-6 1-26 2-3 3-4 4-5 4-7 5-6 5-9 7-8 7-35 8-9 9-37 11-15 12-16
13-14 26-27

G1: [*1-*2], [*3-*4], [*5-*6], [*7-*8]

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G2:N,CH

G3:H,CH3,Et,n-Pr,i-Pr,n-Bu,i-Bu,s-Bu,t-Bu,[*9]

G4:i-Pr,Cy

G5:C,N

Match level :

1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:Atom 8:Atom 9:Atom 10:CLASS
11:CLASS 12:CLASS 13:CLASS 14:CLASS 15:CLASS 16:CLASS 26:CLASS 27:CLASS
32:Atom 35:CLASS 37:CLASS

Generic attributes :

32:

Saturation : Saturated
Type of Ring System : Monocyclic

L10 STRUCTURE UPLOADED

=> d his

(FILE 'REGISTRY' ENTERED AT 17:37:55 ON 04 MAY 2005)

DELETE HISTORY

L1 1163389 S (NC4-C6 OR N2C3-C6 OR NCNC2-C6 OR N3C2-C6)/ES
L2 STRUCTURE UPLOADED
L3 0 S L2 SAMPLE SUB=L1
L4 STRUCTURE UPLOADED
L5 0 S L4 SAMPLE SUB=L1
L6 STRUCTURE UPLOADED
L7 0 S L6 SAMPLE SUB=L1
L8 151 S L6 FULL SUB=L1
L9 1 S C12H16N2 AND L8
L10 STRUCTURE UPLOADED

=> s l10 subset = l8

ENTER SUBSET SEARCH SCOPE - SAMPLE, FULL, RANGE, OR (END):full

FULL SUBSET SEARCH INITIATED 18:30:13 FILE 'REGISTRY'

FULL SUBSET SCREEN SEARCH COMPLETED - 151 TO ITERATE

100.0% PROCESSED 151 ITERATIONS
SEARCH TIME: 00.00.01

24 ANSWERS

L11 24 SEA SUB=L8 SSS FUL L10

=> file caold caplus;

FILE 'CAOLD' ENTERED AT 18:30:35 ON 04 MAY 2005

USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT.

PLEASE SEE "HELP USAGETERMS" FOR DETAILS.

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FILE 'CAPLUS' ENTERED AT 18:30:35 ON 04 MAY 2005

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PLEASE SEE "HELP USAGETERMS" FOR DETAILS.

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=> s l11

10/784,673

L12 2 L11

=> s us-6716837/pn

L13 1 US-6716837/PN

=> s l12 not l13

L14 1 L12 NOT L13

=> d cbib pi hitstr

L14 ANSWER 1 OF 1 CAPLUS COPYRIGHT 2005 ACS on STN

2001:63967 Document No. 134:131423 Preparation of aminoalkylindoles and analogs as 5-HT1D receptor ligands. Edwards, Louise; Isaac, Methvin; Maddaford, Shawn; Slassi, Abdelmalik; Xin, Tao (NPS Allelix Corp., Can.). PCT Int. Appl. WO 2001005758 A2 20010125, 71 pp. DESIGNATED STATES: W: AE, AG, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BY, BZ, CA, CH, CN, CR, CU, CZ, DE, DK, DM, DZ, EE, ES, FI, GB, GD, GE, GH, GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MA, MD, MG, MK, MN, MW, MX, MZ, NO, NZ, PL, PT, RO, RU, SD, SE, SG, SI, SK, SL, TJ, TM, TR, TT, TZ, UA, UG, US, UZ, VN, YU, ZA, ZW, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM; RW: AT, BE, BF, BJ, CF, CG, CH, CI, CM, CY, DE, DK, ES, FI, FR, GA, GB, GR, IE, IT, LU, MC, ML, MR, NE, NL, PT, SE, SN, TD, TG. (English). CODEN: PIXXD2. APPLICATION: WO 2000-CA831 20000714. PRIORITY: US 1999-354091 19990715.

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
PI WO 2001005758	A2	20010125	WO 2000-CA831	20000714
WO 2001005758	A3	20010719		
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RW:	GH, GM, KE, LS, MW, MZ, SD, SL, SZ, TZ, UG, ZW, AT, BE, CH, CY, DE, DK, ES, FI, FR, GB, GR, IE, IT, LU, MC, NL, PT, SE, BF, BJ, CF, CG, CI, CM, GA, GN, GW, ML, MR, NE, SN, TD, TG			
CA 2378047	AA	20010125	CA 2000-2378047	20000714
EP 1196380	A2	20020417	EP 2000-945511	20000714
R:	AT, BE, CH, DE, DK, ES, FR, GB, GR, IT, LI, LU, NL, SE, MC, PT, IE, SI, LT, LV, FI, RO			
JP 2003505369	T2	20030212	JP 2001-511419	20000714

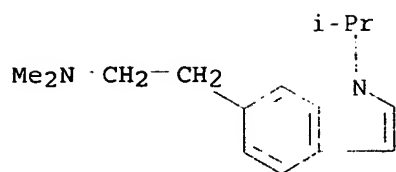
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321745-31-7P 321745-32-8P 321745-33-9P
321745-34-0P 321745-35-1P 321745-37-3P
321745-39-5P 321745-41-9P 321745-45-3P
321745-47-5P 321745-49-7P

RL: BAC (Biological activity or effector, except adverse); BSU (Biological study, unclassified); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)
(preparation of aminoalkylindoles and analogs as 5-HT1D receptor ligands)

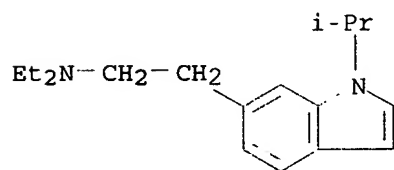
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CN 1H-Indole-6-ethanamine, N,N-dimethyl-1-(1-methylethyl)- (9CI) (CA INDEX NAME)

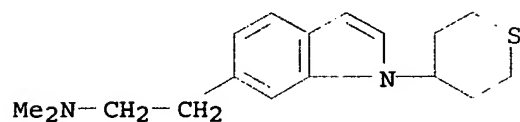
10/784,673



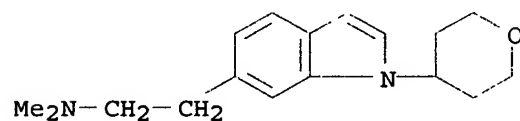
RN 321744-79-0 CAPLUS
CN 1H-Indole-6-ethanamine, N,N-diethyl-1-(1-methylethyl)- (9CI) (CA INDEX NAME)



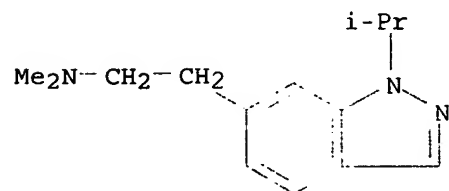
RN 321744-81-4 CAPLUS
CN 1H-Indole-6-ethanamine, N,N-dimethyl-1-(tetrahydro-2H-thiopyran-4-yl)- (9CI) (CA INDEX NAME)



RN 321744-82-5 CAPLUS
CN 1H-Indole-6-ethanamine, N,N-dimethyl-1-(tetrahydro-2H-pyran-4-yl)- (9CI) (CA INDEX NAME)



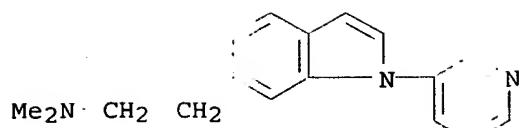
RN 321744-83-6 CAPLUS
CN 1H-Indazole-6-ethanamine, N,N-dimethyl-1-(1-methylethyl)- (9CI) (CA INDEX NAME)



RN 321744-94-9 CAPLUS
CN 1H-Indole-6-ethanamine, N,N-dimethyl-1-(3-pyridinyl)- (9CI) (CA INDEX NAME)

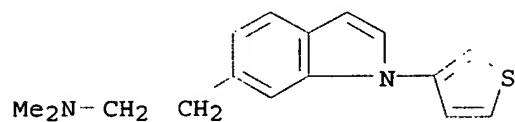
10/784,673

NAME)



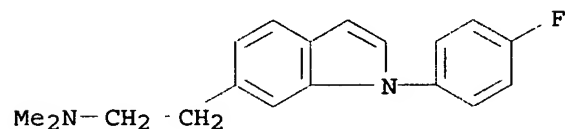
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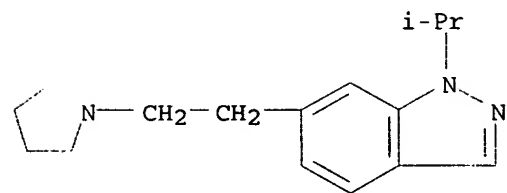
RN 321744-96-1 CAPLUS

CN 1H-Indole-6-ethanamine, 1-(4-fluorophenyl)-N,N-dimethyl- (9CI) (CA INDEX NAME)



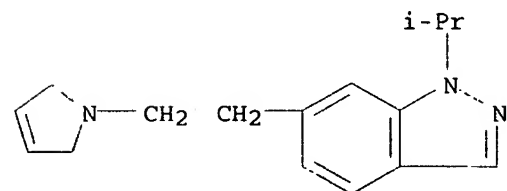
RN 321745-29-3 CAPLUS

CN 1H-Indazole, 1-(1-methylethyl)-6-[2-(1-pyrrolidinyl)ethyl]- (9CI) (CA INDEX NAME)



RN 321745-31-7 CAPLUS

CN 1H-Indazole, 6-[2-(2,5-dihydro-1H-pyrrol-1-yl)ethyl]-1-(1-methylethyl)- (9CI) (CA INDEX NAME)

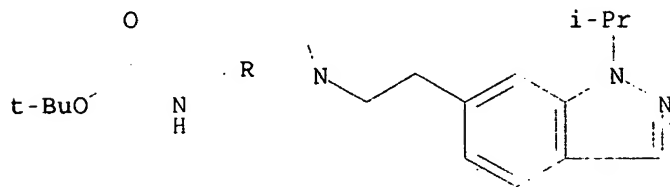


10/784,673

RN 321745-32-8 CAPLUS

CN Carbamic acid, [(3R)-1-[2-[1-(1-methylethyl)-1H-indazol-6-yl]ethyl]-3-pyrrolidinyl]-, 1,1-dimethylethyl ester (9CI) (CA INDEX NAME)

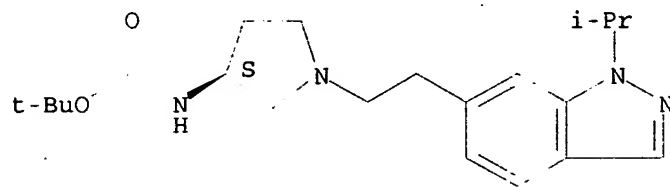
Absolute stereochemistry.



RN 321745-33-9 CAPLUS

CN Carbamic acid, [(3S)-1-[2-[1-(1-methylethyl)-1H-indazol-6-yl]ethyl]-3-pyrrolidinyl]-, 1,1-dimethylethyl ester (9CI) (CA INDEX NAME)

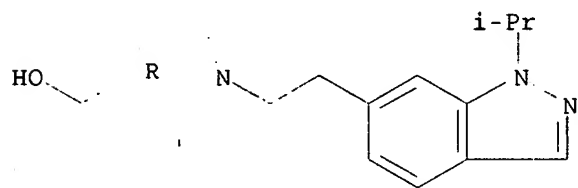
Absolute stereochemistry.



RN 321745-34-0 CAPLUS

CN 3-Pyrrolidinemethanol, 1-[2-[1-(1-methylethyl)-1H-indazol-6-yl]ethyl]-, (3R)- (9CI) (CA INDEX NAME)

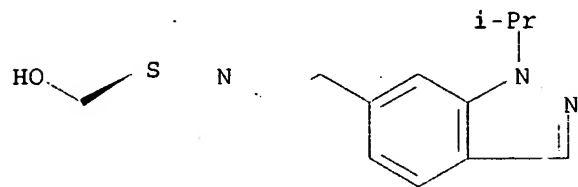
Absolute stereochemistry.



RN 321745-35-1 CAPLUS

CN 3-Pyrrolidinemethanol, 1-[2-[1-(1-methylethyl)-1H-indazol-6-yl]ethyl]-, (3S)- (9CI) (CA INDEX NAME)

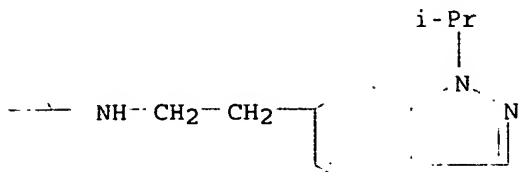
Absolute stereochemistry.



10/784,673

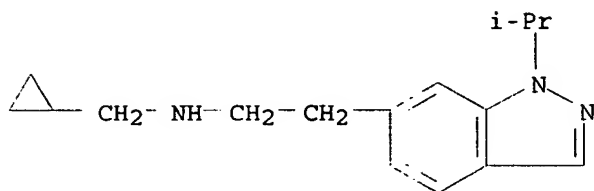
RN 321745-37-3 CAPLUS

CN 1H-Indazole-6-ethanamine, N-cyclopropyl-1-(1-methylethyl)- (9CI) (CA INDEX NAME)



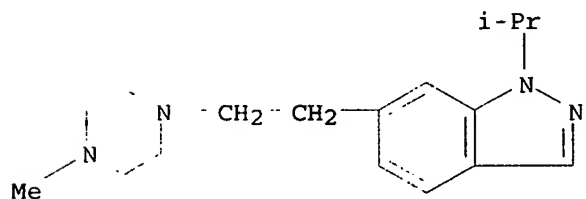
RN 321745-39-5 CAPLUS

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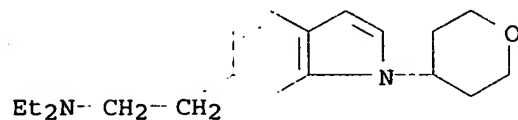
RN 321745-41-9 CAPLUS

CN 1H-Indazole, 1-(1-methylethyl)-6-[2-(4-methyl-1-piperazinyl)ethyl]- (9CI) (CA INDEX NAME)



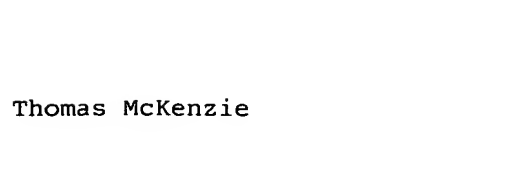
RN 321745-45-3 CAPLUS

CN 1H-Indole-6-ethanamine, N,N-diethyl-1-(tetrahydro-2H-pyran-4-yl)- (9CI) (CA INDEX NAME)

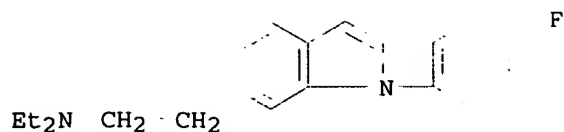


RN 321745-47-5 CAPLUS

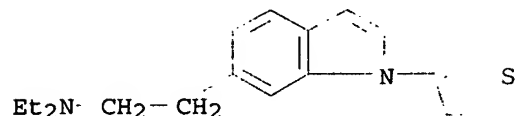
CN 1H-Indole-6-ethanamine, N,N-diethyl-1-(4-fluorophenyl)- (9CI) (CA INDEX NAME)



10/784,673



RN 321745-49-7 CAPLUS
CN 1H-Indole-6-ethanamine, N,N-diethyl-1-(3-thienyl)- (9CI) (CA INDEX NAME)



=> d his

(FILE 'REGISTRY' ENTERED AT 17:37:55 ON 04 MAY 2005)

DELETE HISTORY

L1 1163389 S (NC4-C6 OR N2C3-C6 OR NCNC2-C6 OR N3C2-C6)/ES
L2 STRUCTURE UPLOADED
L3 0 S L2 SAMPLE SUB=L1
L4 STRUCTURE UPLOADED
L5 0 S L4 SAMPLE SUB=L1
L6 STRUCTURE UPLOADED
L7 0 S L6 SAMPLE SUB=L1
L8 151 S L6 FULL SUB=L1
L9 1 S C12H16N2 AND L8
L10 STRUCTURE UPLOADED
L11 24 S L10 SUB=L8 FULL

FILE 'CAOLD, CAPLUS' ENTERED AT 18:30:35 ON 04 MAY 2005

L12 2 S L11
L13 1 S US-6716837/PN
L14 1 S L12 NOT L13

=> s l8

L15 58 L8

=> s l15 not l12

L16 56 L15 NOT L12

=> sort py l16

SORT ENTIRE ANSWER SET? (Y)/N:.

2 ANSWERS DID NOT HAVE 'PY' SORT FIELD

PROCESSING COMPLETED FOR L16

L17 56 SORT L16 PY

=> file reg

FILE 'REGISTRY' ENTERED AT 18:33:27 ON 04 MAY 2005

USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT.

PLEASE SEE "HELP USAGETERMS" FOR DETAILS.

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Property values tagged with IC are from the ZIC/VINITI data file

10/784,673

provided by InfoChem.

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STRUCTURE FILE UPDATES:      3 MAY 2005  HIGHEST RN 849720-40-7
DICTIONARY FILE UPDATES:    3 MAY 2005  HIGHEST RN 849720-40-7
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New CAS Information Use Policies, enter HELP USAGETERMS for details.

TSCA INFORMATION NOW CURRENT THROUGH JANUARY 18, 2005

Please note that search-term pricing does apply when conducting SmartSELECT searches.

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*
* The CA roles and document type information have been removed from
* the IDE default display format and the ED field has been added,
* effective March 20, 2005.  A new display format, IDERL, is now
* available and contains the CA role and document type information.
*
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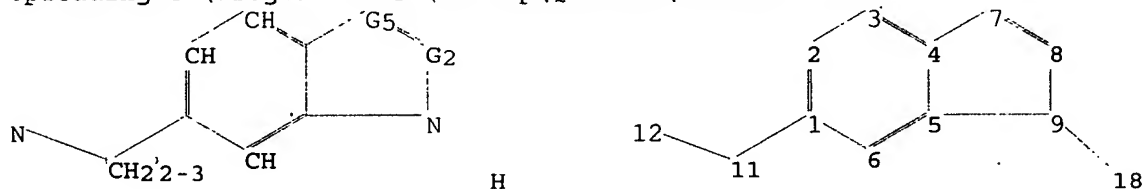
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Crossover limits have been increased. See HELP CROSSOVER for details.

Experimental and calculated property data are now available. For more information enter HELP PROP at an arrow prompt in the file or refer to the file summary sheet on the web at:
<http://www.cas.org/ONLINE/DBSS/registryss.html>

 \Rightarrow

Uploading C:\Program Files\Stnexp\Queries\10784673.str



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chain nodes :
11 18
ring nodes :
1 2 3 4 5 6 7 8 9
ring/chain nodes :
12
chain bonds :
1-11 9-18 11-12
ring bonds :
1-2 1-6 2-3 3-4 4-5 4-7 5-6 5-9 7-8 8-9
exact/norm bonds :
1-11 4-7 5-9 7-8 8-9 9-18 11-12
normalized bonds :
1-2 1-6 2-3 3-4 4-5 5-6
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G1

G2 : N, CH

G5 : C, N

10/784,673

Match level :

1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:Atom 8:Atom 9:Atom 11:CLASS
12:CLASS 18:CLASS

L18 STRUCTURE UPLOADED

=> s l18 subset = l8 full

FULL SUBSET SEARCH INITIATED 18:33:55 FILE 'REGISTRY'

FULL SUBSET SCREEN SEARCH COMPLETED - 151 TO ITERATE

100.0% PROCESSED 151 ITERATIONS

113 ANSWERS

SEARCH TIME: 00.00.01

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L20 38 L8 NOT L19

=> file caold caplus; s l20

FILE 'CAOLD' ENTERED AT 18:34:31 ON 04 MAY 2005

USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT.

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FILE 'CAPLUS' ENTERED AT 18:34:31 ON 04 MAY 2005

USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT.

PLEASE SEE "HELP USAGETERMS" FOR DETAILS.

COPYRIGHT (C) 2005 AMERICAN CHEMICAL SOCIETY (ACS)

L21 8 L20

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L22 6 L21 NOT L12

=> sort py l22

SORT ENTIRE ANSWER SET? (Y)/N:.

PROCESSING COMPLETED FOR L22

L23 6 SORT L22 PY

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L23 ANSWER 1 OF 6 CAPLUS COPYRIGHT 2005 ACS on STN

1976:16500 Document No. 84:16500 Photocyclization of N-chloroacetyl derivatives of indolyethylamines to azepinoindoles and azocinoindoles. Correlation of the reactivity of indole radicals with frontier electron densities calculated by unrestricted Hartree-Fock MO. Naruto, Shunji; Yonemitsu, Osamu (Cent. Res. Lab., Tokyo, Japan). Tetrahedron Letters (39), 3399-402 (English) 1975. CODEN: TELEAY. ISSN: 0040-4039.

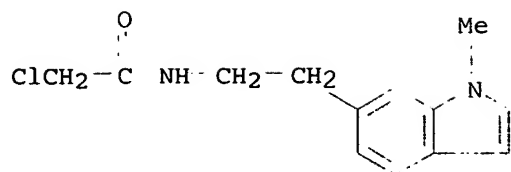
IT 57964-71-3

RL: RCT (Reactant); RACT (Reactant or reagent)

(photochem. cyclization of, MO calcns. for)

RN 57964-71-3 CAPLUS

CN Acetamide, 2-chloro-N-[2-(1-methyl-1H-indol-6-yl)ethyl]- (9CI) (CA INDEX NAME)



L23 ANSWER 2 OF 6 CAPLUS COPYRIGHT 2005 ACS on STN

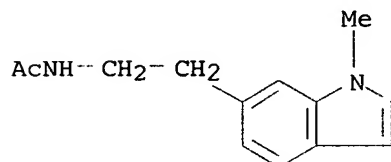
1980:549357 Document No. 93:149357 Photochemical synthesis of azepinoindoles and azocinoindoles from N-chloroacetylinolylethylamines, and a mechanistic study based on the correlation between quantum yields and calculated frontier electron densities of indole radicals. Naruto, Shunji; Yonemitsu, Osamu (Cent. Res. Lab., Sankyo Co., Tokyo, 140, Japan). Chemical & Pharmaceutical Bulletin, 28(3), 900-9 (English) 1980. CODEN: CPBTAL. ISSN: 0009-2363.

IT 74631-90-6P

RL: FORM (Formation, nonpreparative); PREP (Preparation)
(formation of, by photolysis of chloroacetylinolylethylamine)

RN 74631-90-6 CAPLUS

CN Acetamide, N-[2-(1-methyl-1H-indol-6-yl)ethyl]- (9CI) (CA INDEX NAME)



L23 ANSWER 3 OF 6 CAPLUS COPYRIGHT 2005 ACS on STN

1998:42394 Document No. 128:102084 Preparation of 4-heterocyclyl-1-piperidineacetates as glycoprotein IIb/IIIa receptor antagonists. Allen, David George; Eldred, Colin David; Judkins, Brian David; Mitchell, William Leonard; Scopes, David Ian Carter (Glaxo Group Ltd., UK; Allen, David George; Eldred, Colin David; Judkins, Brian David; Mitchell, William Leonard; Scopes, David Ian Carter). PCT Int. Appl. WO 9749698 A1 19971231, 84 pp. DESIGNATED STATES: W: AL, AM, AT, AU, AZ, BA, BB, BG, BR, BY, CA, CH, CN, CU, CZ, DE, DK, EE, ES, FI, GB, GE, GH, HU, IL, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MD, MG, MK, MN, MW, MX, NO, NZ, PL, PT, RO, RU, SD, SE, SG, SI, SK, SL, TJ, TM, TR, TT, UA, UG, US, VZ, VN, YU, ZW, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM; RW: AT, BE, BF, BJ, CF, CG, CH, CI, CM, DE, DK, ES, FI, FR, GA, GB, GR, IE, IT, LU, MC, ML, MR, NE, NL, PT, SE, SN, TD, TG. (English). CODEN: PIXXD2.

APPLICATION: WO 1997-EP3194 19970619. PRIORITY: GB 1996-13017 19960621; GB 1996-13095 19960621; GB 1996-13018 19960621; GB 1996-13026 19960621.

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
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AU 9732610	A1	19980114	AU 1997-32610	19970619
ZA 9705431	A	19981221	ZA 1997-5431	19970619
CN 1222153	A	19990707	CN 1997-195652	19970619

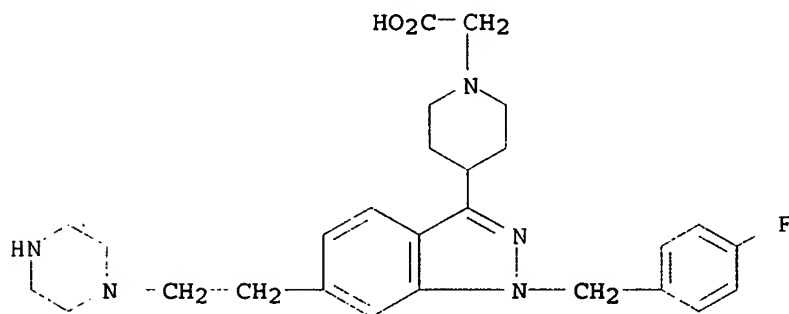
IT 201482-63-5P

RL: BAC (Biological activity or effector, except adverse); BSU (Biological study, unclassified); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)

(preparation of 4-heterocyclyl-1-piperidineacetates as glycoprotein Iib/IIIa receptor antagonists)

RN 201482-63-5 CAPLUS

CN 1-Piperidineacetic acid, 4-[1-[(4-fluorophenyl)methyl]-6-[2-(1-piperazinyl)ethyl]-1H-indazol-3-yl]-, tetrahydrochloride (9CI) (CA INDEX NAME)



● 4 HCl

L23 ANSWER 4 OF 6 CAPLUS COPYRIGHT 2005 ACS on STN

1999:311189 Document No. 130:338114 Indazole bioisostere replacement of catechol in therapeutically active compounds. Marfat, Anthony (Pfizer Products Inc., USA). PCT Int. Appl. WO 9923077 A1 19990514, 231 pp.
 DESIGNATED STATES: W: AL, AM, AT, AU, AZ, BA, BB, BG, BR, BY, CA, CH, CN, CU, CZ, DE, DK, EE, ES, FI, GB, GD, GE, GH, GM, HR, HU, ID, IL, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MD, MG, MK, MN, MW, MX, NO, NZ, PL, PT, RO, RU, SD, SE, SG, SI, SK, SL, TJ, TM, TR, TT, UA, UG, US, UZ, VN, YU, ZW, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM; RW: AT, BE, BF, BJ, CF, CG, CH, CI, CM, CY, DE, DK, ES, FI, FR, GA, GB, GR, IE, IT, LU, MC, ML, MR, NE, NL, PT, SE, SN, TD, TG. (English). CODEN: PIXXD2.
 APPLICATION: WO 1998-IB1710 19981026. PRIORITY: US 1997-64229 19971104; US 1997-64187 19971104; US 1997-64024 19971104; US 1997-64228 19971104; US 1997-64198 19971104.

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
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RW:	GH, GM, KE, LS, MW, SD, SZ, UG, ZW, AT, BE, CH, CY, DE, DK, ES, FI, FR, GB, GR, IE, IT, LU, MC, NL, PT, SE, BF, BJ, CF, CG, CI, CM, GA, GN, GW, ML, MR, NE, SN, TD, TG			
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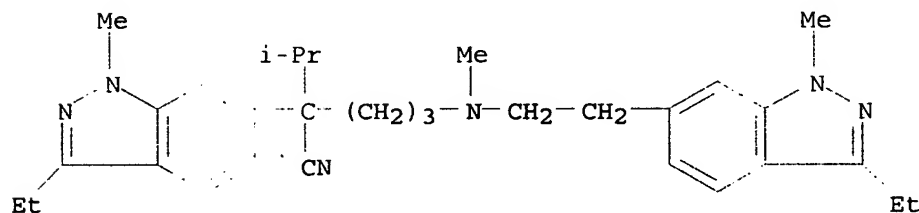
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AU 754734	B2	20021121		
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EP 1028946	A1	20000823	EP 1998-947732	19981026
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JP 2001521926	T2	20011113	JP 2000-518952	19981026
NZ 503918	A	20020328	NZ 1998-503918	19981026
ZA 9810041	A	20000503	ZA 1998-10041	19981103
AP 910	A	20001205	AP 1998-1375	19981104
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US 6329412	B1	20011211	US 2000-535359	20000324
NO 2000002129	A	20000703	NO 2000-2129	20000426
HR 2000000253	A1	20010630	HR 2000-253	20000428
BG 104450	A	20001229	BG 2000-104450	20000517

IT 224188-05-0P

RL: BAC (Biological activity or effector, except adverse); BSU (Biological study, unclassified); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)
(indazole bioisostere replacement of catechol in therapeutically active compds.)

RN 224188-05-0 CAPLUS

CN 1H-Indazole-6-acetonitrile, 3-ethyl- α -[3-[[2-(3-ethyl-1-methyl-1H-indazol-6-yl)ethyl]methylamino]propyl]-1-methyl- α -(1-methylethyl)-(9CI) (CA INDEX NAME)



L23 ANSWER 5 OF 6 CAPLUS COPYRIGHT 2005 ACS on STN

2000:68450 Document No. 132:107953 Preparation of heterocyclic compounds as antagonists of gonadotropin releasing hormone. Goulet, Mark; Ashton, Wallace T.; Chu, Lin; Fisher, Michael H.; Lin, Peter; Ponpipom, Mitree M.; Wyvratt, Matthew J.; Girotra, Narindar N.; Young, Jonathan (Merck and Co., Inc., USA). PCT Int. Appl. WO 2000004013 A1 20000127, 138 pp. DESIGNATED STATES: W: AE, AL, AM, AU, AZ, BA, BB, BG, BR, BY, CA, CN, CU, CZ, EE, GD, GE, HR, HU, ID, IL, IN, IS, JP, KG, KR, KZ, LC, LK, LR, LT, LV, MD, MG, MK, MN, MX, NO, NZ, PL, RO, RU, SG, SI, SK, TJ, TM, TR, TT, UA, US, UZ, VN, YU, ZA, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM; RW: AT, BE, BF, BJ, CF, CG, CH, CI, CM, CY, DE, DK, ES, FI, FR, GA, GB, GR, IE, IT, LU, MC, ML, MR, NE, NL, PT, SE, SN, TD, TG. (English). CODEN: PIXXD2.
APPLICATION: WO 1999-US15581 19990709. PRIORITY: US 1998-115497 19980714.
PATENT NO. KIND DATE APPLICATION NO. DATE

PI	WO 2000004013	A1	20000127	WO 1999-US15581	19990709
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RW: GH, GM, KE, LS, MW, SD, SL, SZ, UG, ZW, AT, BE, CH, CY, DE, DK, ES, FI, FR, GB, GR, IE, IT, LU, MC, NL, PT, SE, BF, BJ, CF, CG, CI, CM, GA, GN, GW, ML, MR, NE, SN, TD, TG

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AU 760030	B2	20030508		
EP 1095038	A1	20010502	EP 1999-933850	19990709
EP 1095038	B1	20030903		

R: AT, BE, CH, DE, DK, ES, FR, GB, GR, IT, LI, LU, NL, SE, MC, PT, IE, SI, LT, LV, FI, RO

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IT 255863-35-5P

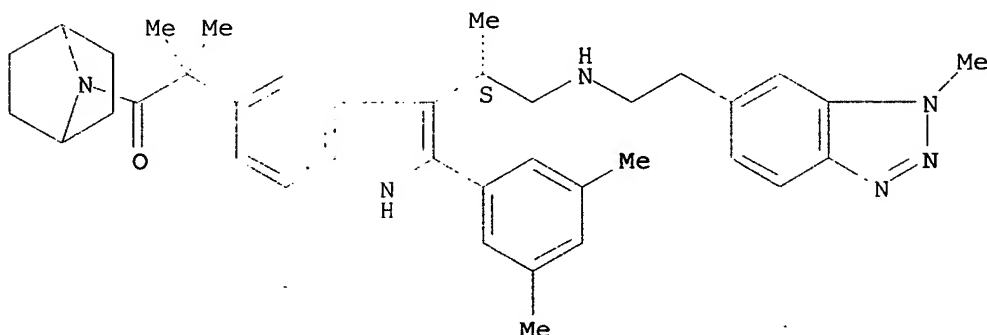
RL: BAC (Biological activity or effector, except adverse); BSU (Biological study, unclassified); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)

(preparation of heterocyclic compds. as antagonists of gonadotropin releasing hormone)

RN 255863-35-5 CAPLUS

CN 7-Azabicyclo[2.2.1]heptane, 7-[2-[2-(3,5-dimethylphenyl)-3-[(1S)-1-methyl-2-[[2-(1-methyl-1H-benzotriazol-6-yl)ethyl]amino]ethyl]-1H-indol-5-yl]-2-methyl-1-oxopropyl]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.



L23 ANSWER 6 OF 6 CAPLUS COPYRIGHT 2005 ACS on STN

2001:178434 Document No. 134:222629 Preparation of indoles as antagonists of gonadotropin releasing hormone. Goulet, Mark; Ashton, Wallace T.; Chu, Lin; Fisher, Michael H.; Lin, Peter; Ponpipom, Mitree M.; Wyvratt, Matthew J.; Girotra, Narindar N.; Young, Jonathan (Merck & Co., Inc., USA). U.S. US 6200957 B1 20010313, 53 pp., Cont.-in-part of U.S. 5,780,437. (English). CODEN: USXXAM. APPLICATION: US 1998-115497 19980714.

PRIORITY: US 1995-PV8633 19951214; US 1996-760816 19961205.

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	JP 2001106685	A2	20010417	JP 2000-257791	19961210
	ZA 9610536	A	19970814	ZA 1996-10536	19961213
	CA 2337407	AA	20000127	CA 1999-2337407	19990709
	WO 2000004013	A1	20000127	WO 1999-US15581	19990709
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 RW: GH, GM, KE, LS, MW, SD, SL, SZ, UG, ZW, AT, BE, CH, CY, DE, DK,
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 CI, CM, GA, GN, GW, ML, MR, NE, SN, TD, TG

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AU 760030 B2 20030508

EP 1095038 A1 20010502 EP 1999-933850 19990709

EP 1095038 B1 20030903

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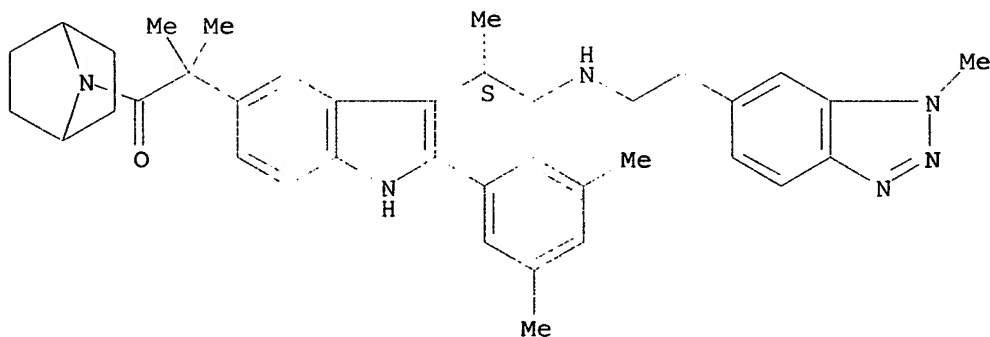
IT 255863-35-5P

RL: BAC (Biological activity or effector, except adverse); BSU (Biological study, unclassified); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)
 (preparation of indoles as antagonists of gonadotropin releasing hormone)

RN 255863-35-5 CAPLUS

CN 7-Azabicyclo[2.2.1]heptane, 7-[2-[2-(3,5-dimethylphenyl)-3-[(1S)-1-methyl-2-[[2-(1-methyl-1H-benzotriazol-6-yl)ethyl]amino]ethyl]-1H-indol-5-yl]-2-methyl-1-oxopropyl]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.



=> d 1 4 cbib pi hitstr

L23 ANSWER 1 OF 6 CAPLUS COPYRIGHT 2005 ACS on STN

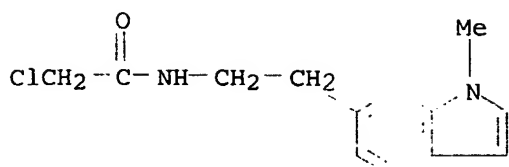
1976:16500 Document No. 84:16500 Photocyclization of N-chloroacetyl derivatives of indolyethylamines to azepinoindoles and azocinoindoles. Correlation of the reactivity of indole radicals with frontier electron densities calculated by unrestricted Hartree-Fock MO. Naruto, Shunji; Yonemitsu, Osamu (Cent. Res. Lab., Tokyo, Japan). Tetrahedron Letters (39), 3399-402 (English) 1975. CODEN: TELEAY. ISSN: 0040-4039.

IT 57964-71-3

RL: RCT (Reactant); RACT (Reactant or reagent)
 (photochem. cyclization of, MO calcns. for)

RN 57964-71-3 CAPLUS

CN Acetamide, 2-chloro-N-[2-(1-methyl-1H-indol-6-yl)ethyl]- (9CI) (CA INDEX NAME)



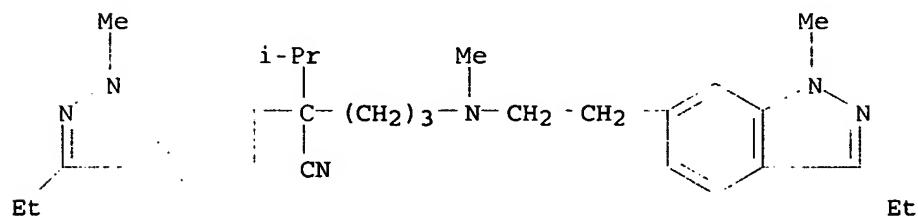
L23 ANSWER 4 OF 6 CAPLUS COPYRIGHT 2005 ACS on STN

1999:311189 Document No. 130:338114 Indazole bioisostere replacement of catechol in therapeutically active compounds. Marfat, Anthony (Pfizer Products Inc., USA). PCT Int. Appl. WO 9923077 A1 19990514, 231 pp.
 DESIGNATED STATES: W: AL, AM, AT, AU, AZ, BA, BB, BG, BR, BY, CA, CH, CN, CU, CZ, DE, DK, EE, ES, FI, GB, GD, GE, GH, GM, HR, HU, ID, IL, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MD, MG, MK, MN, MW, MX, NO, NZ, PL, PT, RO, RU, SD, SE, SG, SI, SK, SL, TJ, TM, TR, TT, UA, UG, US, UZ, VN, YU, ZW, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM; RW: AT, BE, BF, BJ, CF, CG, CH, CI, CM, CY, DE, DK, ES, FI, FR, GA, GB, GR, IE, IT, LU, MC, ML, MR, NE, NL, PT, SE, SN, TD, TG. (English). CODEN: PIXXD2.
 APPLICATION: WO 1998-IB1710 19981026. PRIORITY: US 1997-64229 19971104; US 1997-64187 19971104; US 1997-64024 19971104; US 1997-64228 19971104; US 1997-64198 19971104.

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
PI WO 9923077	A1	19990514	WO 1998-IB1710	19981026
W: AL, AM, AT, AU, AZ, BA, BB, BG, BR, BY, CA, CH, CN, CU, CZ, DE, DK, EE, ES, FI, GB, GD, GE, GH, GM, HR, HU, ID, IL, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MD, MG, MK, MN, MW, MX, NO, NZ, PL, PT, RO, RU, SD, SE, SG, SI, SK, SL, TJ, TM, TR, TT, UA, UG, US, UZ, VN, YU, ZW, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM				
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AU 9894552	A1	19990524	AU 1998-94552	19981026
AU 754734	B2	20021121		
TR 200001234	T2	20000821	TR 2000-200001234	19981026
EP 1028946	A1	20000823	EP 1998-947732	19981026
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BR 9813926	A	20000919	BR 1998-13926	19981026
JP 2001521926	T2	20011113	JP 2000-518952	19981026
NZ 503918	A	20020328	NZ 1998-503918	19981026
ZA 9810041	A	20000503	ZA 1998-10041	19981103
AP 910	A	20001205	AP 1998-1375	19981104
W: BW, GM, GH, KE, MW, SD, UG, ZM, ZW				
US 6391872	B1	20020521	US 1999-381425	19990920
US 6329412	B1	20011211	US 2000-535359	20000324
NO 2000002129	A	20000703	NO 2000-2129	20000426
HR 2000000253	A1	20010630	HR 2000-253	20000428
BG 104450	A	20001229	BG 2000-104450	20000517
IT 224188-05-0P 224188-16-3P				
RL: BAC (Biological activity or effector, except adverse); BSU (Biological study, unclassified); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)				
(indazole bioisostere replacement of catechol in therapeutically active compds.)				
RN 224188-05-0 CAPLUS				

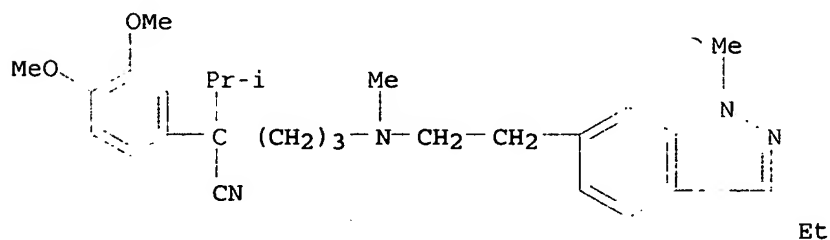
10/784,673

CN 1H-Indazole-6-acetonitrile, 3-ethyl- α -[3-[[2-(3-ethyl-1-methyl-1H-indazol-6-yl)ethyl]methylamino]propyl]-1-methyl- α -(1-methylethyl)-(9CI) (CA INDEX NAME)



RN 224188-16-3 CAPLUS

CN Benzeneacetonitrile, α -[3-[[2-(3-ethyl-1-methyl-1H-indazol-6-yl)ethyl]methylamino]propyl]-3,4-dimethoxy- α -(1-methylethyl)-(9CI) (CA INDEX NAME)



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ALL L# QUERIES AND ANSWER SETS ARE DELETED AT LOGOFF

LOGOFF? (Y)/N/HOLD:.

STN INTERNATIONAL LOGOFF AT 18:39:18 ON 04 MAY 2005

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